Documentation of ModelSolver * A Python class for analyzing dynamic algebraic models[†]

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Abstract

This paper documents the Python class ModelSolver. ModelSolver lets the user define a model object in terms of equations and endogenous variables. It contains methods to solve the model subject to data (in a Pandas DataFrame), as well as analyzing the model using graph theory and network plots.

What sets ModelSolver apart from other similar solvers, is that it does not require the equations of the model to be written in any particular form, or that the user associates equations with endogenous variables. Most other solvers require that either 1) the model is normalized (i.e., that the model is written in terms of endogenous variables), or 2) that the user explicitly associates equations with endogenous variables. This is non-trivial for models with lots of equations. ModelSolver, however, reads equations in whatever form they may be formulated, and performs the necessary analyzes, without any input from the user other than lists of equations and endogenous variables.

ModelSolver was developed in order to facilitate solving an inputoutput model for the Norwegian monthly national accounts. It analyzes and solves the model's more than 15,500 equations over more than 30 periods in under a minute on a laptop computer.

^{*}Code in https://github.com/statisticsnorway/model-solver.git. Repository also contains Notebook with examples of applications.

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1 Background

1.1 The model

Consider a model consisting of n equations, ¹

$$L_1(\mathbf{x}_t|\mathbf{z}_t) = R_1(\mathbf{x}_t|\mathbf{z}_t),$$

$$L_2(\mathbf{x}_t|\mathbf{z}_t) = R_2(\mathbf{x}_t|\mathbf{z}_t),$$

$$\vdots$$

$$L_n(\mathbf{x}_t|\mathbf{z}_t) = R_n(\mathbf{x}_t|\mathbf{z}_t),$$

for $t = T_0, T_0 + 1, ..., T_1$, where $\mathbf{x}_t = (x_{1,t}, x_{2,t}, ..., x_{n,t})$ is a vector of endogenous variables (and, the way it is written, the values for which solve the model), and $\mathbf{z}_t = (z_{1,t-k_1}, z_{2,t-k_2}, ..., z_{m,t-k_m}, ..., x_{i,t-\kappa_i}, ...)$, for $k_i \geq 0$ and $\kappa_i > 0$, is a vector of contemporaneous and lags of exogenous variables and lags of endogenous variables (or taken together: predetermined variables).
² It's convenient to formulate the model using a vector function $\mathbf{F} : \mathbb{R}^n \mapsto \mathbb{R}^n$,
³

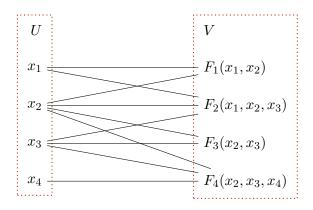
$$\mathbf{F}(\mathbf{x}_t|\mathbf{z}_t) = \begin{pmatrix} F_1(\mathbf{x}_t|\mathbf{z}_t) \\ F_2(\mathbf{x}_t|\mathbf{z}_t) \\ \vdots \\ F_n(\mathbf{x}_t|\mathbf{z}_t) \end{pmatrix} = \begin{pmatrix} L_1(\mathbf{x}_t|\mathbf{z}_t) - R_1(\mathbf{x}_t|\mathbf{z}_t) \\ L_2(\mathbf{x}_t|\mathbf{z}_t) - R_2(\mathbf{x}_t|\mathbf{z}_t) \\ \vdots \\ L_n(\mathbf{x}_t|\mathbf{z}_t) - R_n(\mathbf{x}_t|\mathbf{z}_t) \end{pmatrix},$$

 $^{^{1}}L_{i}$ and R_{i} denote the left and right hand side of equation i, respectively. L_{i} or R_{i} may be as simple as $x_{j,t}$ or a constant, or something more complicated involving many endogenous variables.

²Note that \mathbf{z}_t may contain only $z_{i,t-k_i}$'s or $x_{i,t-k_i}$'s or be empty.

 $^{^3}$ We say that **F** maps from \mathbb{R}^n because exogenous variables (contemporaneous and lags) and lags of the endogenous variables are taken as given in any given period, i.e., predetermined.

Figure 1: Bipartite graph (BiGraph of model)



with the solution to the model being given by $\{\mathbf{x}_t\}_{t=T_0}^{T_1}$ such that

$$\mathbf{F}(\mathbf{x}_t|\mathbf{z}_t) = \mathbf{0} \text{ for } t = T_0, T_0 + 1, \dots, T_1.$$

On the face of it, this is a problem which the Newton-Raphson-algorithm handles well. The issue, however, is that n might be quite large. In the Norwegian national accounts, for instance, n is greater than 15,500. Therefore, it is useful to analyze the system of equations before solving it, in order to break it down into minimal $simultaneous\ blocks$ that can be solved in a particular order.

1.2 Block analysis and ordering

In order to analyze and divide the model into blocks, we use results from graph theory.

First, we construct a bipartite graph (BiGraph) that connects endogenous variables with equations; i.e., B = (U, V, E), where U is a set containing nodes representing endogenous varibales, V is a set containing nodes representing equations, and E is a set containing edges between nodes in U and V. This is illustrated in Figure 1 for an arbitrary model with four equations (we omit time subscripts and exogenous variables and lags for notational convenience).

Next, we apply maximum bipartite matching (MBM), which assigns each endogenous variable to one and only one equation. Formally, MBM results in a function or mapping M, with M(x) = F and $M^{-1}(F) = x$ for $x \in U$ and $F \in V$ (i.e., M is a one-to-one mapping from endogenous variables to equations, $M: U \mapsto V$ and $M^{-1}: V \mapsto U$). The MBM is arbitrary (G may have as many as n! MBMs), and Figure 2 shows one possible solution for our arbitrary model.

We use the MBM to determine what endogenous variables impact what *other* endogenous variables. We do this by 1) assigning endogenous variables

Figure 2: Maximum Bipartite Match (MBM) of BiGraph

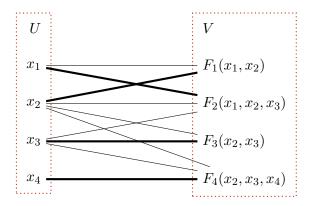
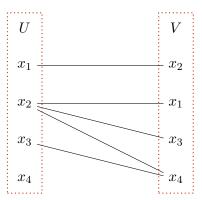


Figure 3: Graph of what endogenous variables impact what *other* endogenous variables

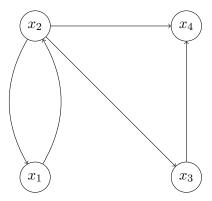


to equations (or 'labeling' the equations 'as' endogenous variables) using $M^{-1}(F) = x$ (the MBM, after all, is a one-to-one mapping between endogenous variables and equations), and 2) removing the MBM edges (the MBM edges simply imply that each endogenous variable impacts itself, which, although not altering the end result, is non-informative). This procedure leaves us with a new BiGraph B' = (U, V', E'), where the nodes in V' are the re-labeled nodes in V (i.e., $V' = \{M^{-1}(F)|F \in V\}$), and E' contains all the edges except the MBM edges. This is illustrated in Figure 3 for our arbitrary model.

Note that, although the MBM is arbitrary, the description of what endogenous variables affect what *other* endogenous variables E' is *not* (see proof).

Next, we use B' to construct a directed graph (DiGraph) that shows how the endogenous variables impact each other (the causality goes from U to V in B'); i.e., D = (V, A), where V contains nodes representing endogenous

Figure 4: Directed graph (DiGraph) of what endogenous variables impact what *other* endogenous variables



variables, and A contains directed edges or arcs from endogenous variables to other endogenous variables. The DiGraph is shown in Figure 4.

Finally, we find the *strong components* of the D, as shown in Figure 5 for our arbitrary model. A strong component is a set of nodes that are connected such that every node can be reached from every other node in the set (traversing the arcs). A *condensation* of the DiGraph is a (new) DiGraph with each node being a strong component (of the former). The condensation can be illustrated as in Figure 6 for our arbitrary model. Each node of the condensation corresponds to a block of the model, and the arcs decide the order in which the blocks are to be solved. See [1] for a discussion of strong components and block analysis.

Note that, since the MBM is a one-to-one-mapping, we may draw the DiGraph and condensation labeling the nodes with endogenous variables *or* equations, switching back and forth using the MBM.

In our arbitrary model, x_1 and x_2 must be solved first in one simultaneous block. Next, x_3 is solved (taking x_1 and x_2 as given, being predetermined by the solution to block 1). Finally, x_4 is solved (taking x_1 , x_2 and x_3 as given, being predetermined by the solutions to block 1 and 2).

1.3 Simulation code

With the blocks of the model as given by the condensation of the model DiGraph, as previously discussed, we can generate *simulation code* for each block. That is *either*

Figure 5: Strong components of DiGraph

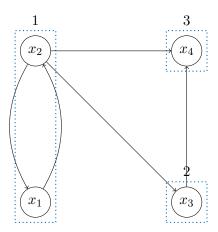
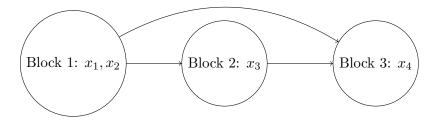


Figure 6: Condensed DiGraph



- a symbolic function (definition) that takes predetermined input and returns the endogenous value (if the block is a definition, as described below), or
- a symbolic objective function and Jacobian matrix that take predetermined input and (initial) values for the endogenous variables, which are to be sent to a Newton-Raphson algorithm, which in turn returns the endogenous value(s)

A block is a said to be a definition if and only if it 1) consists of one equation and 2) the only thing on the left hand side of the equation is the endogenous variable of that equation (and that endogenous variable does *not* show up on the right hand side of that equation too).

1.4 Solution

In order to solve the model numerically, we loop (nested) over time periods (chronologically) and blocks (according to the order dictated by the condensation of the model DiGraph).

If the block to be solved is a definition, as discussed above, the solution is given straight forwardly by the function defining it (given predetermined input).

If the block to be solved is not a definition, it is sent to a Newton-Raphson algorithm, which in turn returns the solution (given predetermined input, and initial values for the endogenous variables).

The k + 1st iteration of the Newton-Raphson algorithm is given by

$$\mathbf{x}_t^{(k+1)} = \mathbf{x}_t^{(k)} - \mathbf{J}_{\mathbf{F}}^{-1}(\mathbf{x}_t^{(k)}|\mathbf{z}_t)\mathbf{F}(\mathbf{x}_t^{(k)}|\mathbf{z}_t),$$

where $\mathbf{J}_{\mathbf{F}}(\mathbf{x}_t|\mathbf{z}_t)$ is the Jacobian matrix of $\mathbf{F}(\mathbf{x}_t|\mathbf{z}_t)$ evaluated at \mathbf{x}_t and \mathbf{z}_t . The algorithm stops and returns $\mathbf{x}_t^{(k+1)}$ if $\max\left(\left|\mathbf{x}_t^{(k+1)}-\mathbf{x}_t^{(k)}\right|\right) \leq \varepsilon$, where ε is a tolerance level. If it so happens that $\mathbf{F}(\mathbf{x}_t^{(0)}|\mathbf{z}_t) = \mathbf{0}$, then the initial values solve the model, and as such $\mathbf{x}_t^{(0)}$ is returned as the solution.

2 Examples of use

The ModelSolver object is instantiated by

model = ModelSolver(equations, endogenous)

where equations and endogenous are lists containing equations and endogenous variables as strings, e.g.,

```
equations = [
'x1 = a1',
'x2 = a2',
'0.2*x1+0.7*x2 = 0.1*ca+0.8*cb+0.3*i1',
'0.8*x1+0.3*x2 = 0.9*ca+0.2*cb+0.1*i2',
'k1 = k1(-1)+i1',
'k2 = k2(-1)+i2'
]
endogenous = ['x1', 'x2', 'ca', 'cb', 'k1', 'k2']
```

Note that the (-#)-syntax denotes lags of a variable; ModelSolver does not support leads. Mathematical functions supported are max(), min(), log() and exp(). Note also that the equations are written in a form that is neither normalized, nor associates equations with endogenous variables.

It is possible to switch endogenous and exogenous variables using

```
model.switch_endo_vars(old_endo_vars, new_endo_vars)
```

where old_endo_vars and new_endo_vars are lists. Upon invoking switch_endo_vars, a tuple of endogenous variables stored inside the model object is created, and ModelSolver immediately performs block analysis on the updated model, such that it can be solved for the updated list of endogenous variables.

2.1 Solving the model

Let input_df be a Pandas DataFrame containing exogenous variables and initial values for the endogenous variables. Then

```
solution_df = model.solve(input_df)
```

solves the model and stores the results in solution_df, which is a Pandas DataFrame *identical* to input_df, *but* where the endogenous variables are the solutions to the model.

2.2 Analyzing the model

ModelSolver contains a number of methods that let the user analyze the model:

model.describe() describes the model in terms of number of equations and blocks, along with information about the blocks, e.g.,

Model consists of 6 equations in 5 blocks 4 of the blocks consist of simple definitions

```
4 blocks have 1 equations 1 blocks have 2 equations
```

model.find_endo_var(<endogenous variable>) returns the block in which the endogenous variable is solved for.

model.show_block(<block number>) returns information about the block—endogenous variables, exogenous and predetermined variables and equations—e.g.,

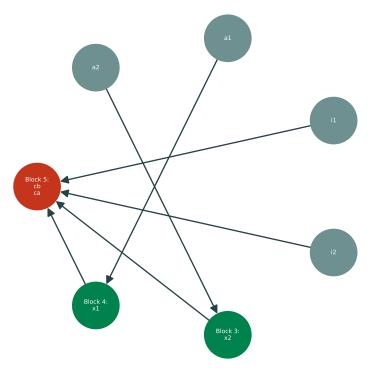
Block consists of an equation or a system of equations

```
2 endogenous variables:
cb ca
4 exogenous or predetermined variables:
i2 i1 x1 x2
2 equations:
0.8*x1+0.3*x2 = 0.9*ca+0.2*cb+0.1*i2
0.2*x1+0.7*x2 = 0.1*ca+0.8*cb+0.3*i1
```

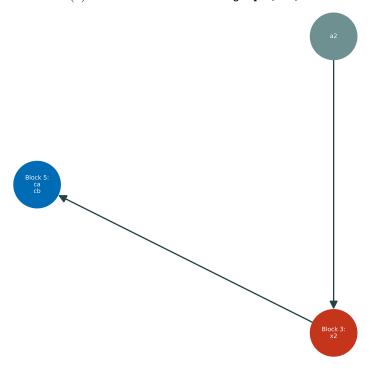
model.show_blocks() returns information about all the blocks of the model.
model.trace_to_exog_vars(<block number>) traces the model DiGraph
from the block back to the nodes of origin and reports what exogenous
variables determine those nodes.

model.trace_to_exog_vals(<block number>, <period index>) traces the model DiGraph back from the block to the nodes of origin, finds what exogenous variables determine those nodes, and reports values for those variables for the chosen period index (0,1,...).

model.draw_blockwise_graph() draws a network graph as shown in Figure 7. The method takes as arguments a variable name, and the maximum number of ancestor and descendant nodes, and maximum number of nodes in total.



 $(a) \ {\tt model.draw_blockwise_graph('ca')}$



(b) model.draw_blockwise_graph('x2')

Figure 7: Examples of network plots

3 Dependencies

ModelSolver is built using the following packages:

- NumPy
- NetworkX
- Pandas
- SymEngine
- Numba
- collections
- functools
- Matplotlib

References

[1] Gilli, Manfred. "Causal Ordering and Beyond." International Economic Review, vol. 33, no. 4, 1992, pp. 957–71. JSTOR, https://doi.org/10.2307/2527152. Accessed 22 Mar. 2023.

A Proofs

Theorem 1. All MBMs M_1, M_2, \ldots result in the same description of what endogenous variables impact what other endogenous variables $E'_1 = E'_2 = \ldots = E'$.

Proof. Suppose we have a BiGraph B = (U, V, E), with

$$E = \{(x_i, F_p), (x_i, F_q), (x_i, F_p), (x_i, F_q)\} \cup \tilde{E},$$

where \tilde{E} is a set containing any *other* edges between endogenous variables and equations (or functions) F than the ones explicitly given. Suppose further that B has two MBMs (B may have more MBMs—in fact as many as n!—but we shall only concern ourselves with these two for the purpose of this proof), ⁴

$$M_1 = \{(x_i, F_p), (x_j, F_q)\} \cup \tilde{M},$$

 $M_2 = \{(x_i, F_q), (x_j, F_p)\} \cup \tilde{M},$

that is, M_1 and M_2 are identical, except for two matches (the ones associating x_i and x_j with F_p and F_q ; \tilde{M} is a set containing all other matches). We use the MBMs to 'label' equations F 'as' endogenous variables, $M^{-1}(F) = x$, as discussed in the main text, i.e.,

$$E'_{1} = \{(x_{i}, x_{i}), (x_{i}, x_{j}), (x_{j}, x_{i}), (x_{j}, x_{j})\} \cup \tilde{E}',$$

$$E'_{2} = \{(x_{i}, x_{j}), (x_{i}, x_{i}), (x_{j}, x_{j}), (x_{j}, x_{i})\} \cup \tilde{E}'.$$

where prime denotes that the set contains equations 'labeled as' endogenous variables using the MBM. (We ignore the mappings from an endogenous variable onto itself.) Note that \tilde{E}' is the same for E'_1 and E'_2 . This follows because we have chosen \tilde{M} to be the same for M_1 and M_2 . Therefore, we see that $E'_1 = E'_2$. Since we can go from any MBM to any other switching two and two edges repeatedly, it follows that any MBM M_1, M_2, \ldots gives the same result as any other $E'_1 = E'_2 = \ldots = E'$, which is what we set out to prove.

⁴As we can choose x_i , x_j , F_p and F_q , this is always the case, unless the MBM is unique (which would render the proof superfluous).